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## IN THE UNITED STATES PATENT AND TRADEMARK OFFICE

IN RE APPLICATION OF

Art Unit: 1625

ROBL ET AL.

Examiner: Evelyn Mei Huang

APPLICATION NO: DIVISION OF APPLICATION SERIAL

NO. 10/008,154 FILED DECEMBER 4, 2001

FILED: HEREWITH

FOR: HMG-COA REDUCTASE INHIBITORS AND METHOD

Commissioner for Patents P.O. Box 1450 Alexandria, VA 22313-1450

## PRE-EXAMINATION AMENDMENT

Sir:

Please amend the claims of the subject application to read as follows.

Please cancel Claims 1 to 11, 13, 16, 37 to 40 and 45.

17. (Amended) A pharmaceutical combination comprising the HMG CoA reductase inhibitor compound having the structure

wherein

Z is 
$$^{HO}$$
  $^{CO_2R_3}$  or  $^{OH}$  also referred to as the  $\delta$ -lactone;

n is 0 or 1;

x is 0, 1, 2, 3 or 4;

y is 0, 1, 2, 3 or 4, provided that at least one of x and y is other than 0; and optionally one or more carbons of  $(CH_2)_x$  and/or one or more carbons of  $(CH_2)_y$  together with additional carbons form a 3 to 7 membered spirocyclic ring;

 $R_1$  and  $R_2$  are the same or different and are independently selected from alkyl, arylalkyl, cycloalkyl, alkenyl, cycloalkenyl, aryl, heteroaryl or cycloheteroalkyl;

R<sub>3</sub> is H or lower alkyl;

R<sub>4</sub> is halogen, CF<sub>3</sub>, hydroxy, alkyl, alkoxy, carboxyl, carboxyalkyl-, aminoalkyl, amino, alkanoylamino, aroylamino, cyano, alkoxyCON(R<sub>10</sub>)-, R<sub>11</sub>R<sub>12</sub>NCO<sub>2</sub>-, R<sub>11</sub>R<sub>12</sub>NCO-, R<sub>13</sub>SO<sub>2</sub>N(R<sub>10</sub>)-, R<sub>11</sub>R<sub>12</sub>NSO<sub>2</sub>N(R<sub>10</sub>)-, R<sub>13</sub>OCO<sub>2</sub>- or R<sub>13</sub>OCO;

R<sub>13</sub> is alkyl, arylalkyl, cycloalkyl, alkenyl, cycloalkenyl, aryl, heteroaryl or cycloheteroalkyl;

 $R_{11}$  and  $R_{12}$ , and  $R_{10}$  are the same or different and are independently selected from H, alkyl, arylalkyl, cycloalkyl, alkenyl, cycloalkenyl, aryl, heteroaryl or cycloheteroalkyl;

or  $R_{11}$  and  $R_{12}$  may be taken together with the nitrogen to which they are attached to form a stable 3 to 8 membered ring, which, where applicable, includes 1 to 3 heteroatoms in the ring.

R<sub>7</sub> is H or lower alkyl;

and represents a single bond or a double bond (which may be cis or trans);

or a pharmaceutically acceptable salt thereof (when R<sub>3</sub> is H), or an ester thereof and or a stereoisomer thereof, and another therapeutic agent which is one or more hypolipidemic agents or lipid-lowering agents, or lipid agents, or lipid modulating agents, and/or one or more other types of therapeutic agents including antidiabetic agents, anti-obesity agents, antihypertensive agents, platelet aggregation inhibitors, anti-dementia agents, anti-Alzheimer's agents, anti-osteoporosis agents, and/or hormone replacement therapeutic agents, and/or other cardiovascular agents (including anti-anginal agents, anti-arrhythmic agents, anti-atherosclerosis agents, anti-inflammatory agents, anti-arrhitis agents, anti-platelet agents, anti-heart failure agents), anti-cancer agents, anti-infective agents, hormone replacement agents, growth hormone secretagogues, selective androgen receptor modulators, and/or immunomodulatory agents.

42. (Amended) A method for treating cholesterol related diseases, diabetes and related diseases, cardiovascular diseases, cerebrovascular diseases, which comprises administering to a mammalian species in need of treatment a therapeutically effective amount of a combination of a compound having the structure

wherein

;)

Z is 
$$R_7$$
 CO<sub>2</sub>R<sub>3</sub> or  $R_7$  also referred to as the  $\delta$ -lactone;

n is 0 or 1;

x is 0, 1, 2, 3 or 4;

y is 0, 1, 2, 3 or 4, provided that at least one of x and y is other than o; and optionally one or more carbons of  $(CH_2)_x$  and/or one or more carbons of  $(CH_2)_y$  together with additional carbons form a 3 to 7 membered spirocyclic ring;

R<sub>1</sub> and R<sub>2</sub> are the same or different and are independently selected from alkyl, arylalkyl, cycloalkyl, alkenyl, cycloalkenyl, aryl, heteroaryl or cycloheteroalkyl;

R<sub>3</sub> is H or lower alkyl;

 $R_4$  is halogen,  $CF_3$ , hydroxy, alkyl, alkoxy, carboxyl, carboxyalkyl-, aminoalkyl, amino, alkanoylamino, aroylamino, cyano, alkoxy $CON(R_{10})$ -,  $R_{11}R_{12}NCO_2$ -,  $R_{11}R_{12}NCO$ -,  $R_{13}SO_2N(R_{10})$ -,  $R_{11}R_{12}NSO_2N(R_{10})$ -,  $R_{13}OCO_2$ - or  $R_{13}OCO_3$ -

R<sub>13</sub> is alkyl, arylalkyl, cycloalkyl, alkenyl, cycloalkenyl, aryl, heteroaryl or cycloheteroalkyl;

R<sub>11</sub> and R<sub>12</sub>, and R<sub>10</sub> are the same or different and are independently selected from H, alkyl, arylalkyl, cycloalkyl, alkenyl, cycloalkenyl, aryl, heteroaryl or cycloheteroalkyl;

R<sub>7</sub> is H or lower alkyl;

and represents a single bond or a double bond (which may be cis or trans);